PROJECTED ANSWERS:

ring bonds :  $1-2 \quad 1-6 \quad 2-3 \quad 3-4 \quad 3-14 \quad 4-5 \quad 4-7 \quad 5-6 \quad 7-8 \quad 7-9 \quad 8-12 \quad 9-10 \quad 9-14 \quad 10-11 \quad 11-12$ exact/norm bonds : 2-15 3-14 4-7 6-18 9-14 normalized bonds : 1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-9 8-12 9-10 10-11 11-12G1:Ir,Rh G2:C1, Br, F, CH3, Et, n-Pr, i-Pr, Ph, MeO, EtO, n-PrO, i-PrO, CN, NO2 Match level : 1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 14:Atom 15:CLASS 18:CLASS L1STRUCTURE UPLOADED => d his (FILE 'HOME' ENTERED AT 10:25:32 ON 22 JUN 2009) FILE 'REGISTRY' ENTERED AT 10:25:54 ON 22 JUN 2009 STRUCTURE UPLOADED L1 => d 11L1 HAS NO ANSWERS L1STR `G2 NО G1 Ir,Rh G2 Cl,Br,F,Me,Et,n-Pr,i-Pr,Ph,MeO,EtO,n-PrO,i-PrO,CN,NO2 Structure attributes must be viewed using STN Express query preparation. => s 11 SAMPLE SEARCH INITIATED 10:26:22 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED -5 TO ITERATE 100.0% PROCESSED 5 ITERATIONS 2 ANSWERS SEARCH TIME: 00.00.01 \*\*COMPLETE\*\* FULL FILE PROJECTIONS: ONLINE BATCH \*\*COMPLETE\*\* PROJECTED ITERATIONS: 5 TO

234

124

2 TO

=> s 11 full

FULL SEARCH INITIATED 10:26:28 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 74 TO ITERATE

100.0% PROCESSED 74 ITERATIONS 9 ANSWERS

SEARCH TIME: 00.00.01

L3 9 SEA SSS FUL L1

=> fil caplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 185.88 186.10

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FILE COVERS 1907 - 22 Jun 2009 VOL 150 ISS 26 FILE LAST UPDATED: 21 Jun 2009 (20090621/ED) REVISED CLASS FIELDS (/NCL) LAST RELOADED: Apr 2009 USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Feb 2009

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 13

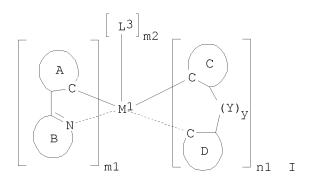
L4 1 L3

=> d bib abs

- L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2009 ACS on STN
- AN 2006:632152 CAPLUS
- DN 145:112952
- TI Metal complexes with nucleophilic carbene ligands and devices and processes using them
- IN Pretot, Roger; Van Der Schaaf, Paul Adriaan; Schmidt, Jemima; Schmidhalter, Beat; Schaefer, Thomas; Lamatsch, Bernd
- PA Ciba Specialty Chemicals Holding Inc., Switz.
- SO PCT Int. Appl., 149 pp. CODEN: PIXXD2

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DT Patent
LA English
FAN.CNT 1
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1 1314 • 4	PATENT NO.				KIND		DATE		APPLICATION NO.						DATE			
ΡI	WO 2006067074			A1		20060629		WO 2005-EP56767						20051214				
	W:	ΑE,	AG,	AL,	AM,	ΑT,	AU,	AZ,	BA,	BB,	BG,	BR,	BW,	BY,	BZ,	CA,	CH,	
		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,	
		GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KM,	KN,	KP,	KR,	
		KZ,	LC,	LK,	LR,	LS,	LT,	LU,	LV,	LY,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	
		MZ,	NA,	NG,	NΙ,	NO,	NZ,	OM,	PG,	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,	
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		VN,	YU,	ZA,	ZM,	ZW												
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		IS,	ΙΤ,	LT,	LU,	LV,	MC,	NL,	PL,	PT,	RO,	SE,	SI,	SK,	TR,	BF,	ВJ,	
		CF,	CG,	CI,	CM,	GΑ,	GN,	GQ,	GW,	ML,	MR,	NE,	SN,	TD,	ΤG,	BW,	GH,	
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		KG,	KΖ,	MD,	RU,	ΤJ,	TM											
	CA 2589711				A1				CA 2005-2589711									
	EP 1841834				A1	1 20071010			EP 2005-817212						20051214			
	EP 1841834			В1		2009	0506											
	R:	ΑT,															ΙE,	
		,	,		•	,	LV,	MC,	NL,	PL,	PT,	RO,	SE,	SI,	SK,	TR		
	CN 101087863								CN 2005-80044163									
	JP 2008525366				Τ	20080717			JP 2007-547457									
	AT 430789								AT 2005-817212									
	IN 2007CN02780								IN 2007-CN2780									
	KR 2007091355								KR 2007-716762						20070720			
PRAI	EP 2004						2004											
	WO 2005-EP56767						2005	1214										
OS	MARPAT	145:	1129	52														
GI																		



а

AB The title complexes are described by the general formula I (ring A = an optionally substituted aryl group which can optionally contain heteroatoms; ring B = an optionally substituted nitrogen-containing aryl group, which can optionally contain further heteroatoms; ring A and ring B may be bonded to form a ring; group C = an acyclic carbene or a cyclic carbene which can optionally contain heteroatoms; ring D = an optionally substituted aryl group which can optionally contain heteroatoms; n1 = 1 = 3, m1 = 0, 1, or 2; m2 = 0 or 1; M1 = a metal with an atomic weight > 40; L3 =

monodentate or bidentate ligand; Y = -C(=0) - or -C(X1)2-; X1 = H or C1-4 alkyl; and y = 0 or 1) with the exception of certain specified compds. The use of the compds. is described in electronic devices, especially organic

light-emitting diodes, as oxygen-sensitive indicators, as phosphorescent indicators in bioassays, and as catalysts. Organic electronic devices, especially

organic light-emitting diodes, comprising an emitting layer which comprises the compds., as well as displays employing the light-emitting diodes, are also described.

RE.CNT 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

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=>

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=> LOG Y

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NEWS 9 APR 28 Limits doubled for structure searching in CAS REGISTRY

- NEWS 10 MAY 08 STN Express, Version 8.4, now available
- NEWS 11 MAY 11 STN on the Web enhanced
- NEWS 12 MAY 11 BEILSTEIN substance information now available on STN Easy
- NEWS 13 MAY 14 DGENE, PCTGEN and USGENE enhanced with increased limits for exact sequence match searches and introduction of free HIT display format
- NEWS 14 MAY 15 INPADOCDB and INPAFAMDB enhanced with Chinese legal status data
- NEWS 15 MAY 28 CAS databases on STN enhanced with NANO super role in records back to 1992
- NEWS 16 JUN 01 CAS REGISTRY Source of Registration (SR) searching enhanced on STN

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STRUCTURE FILE UPDATES: 21 JUN 2009 HIGHEST RN 1159253-26-5 DICTIONARY FILE UPDATES: 21 JUN 2009 HIGHEST RN 1159253-26-5

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=>

Uploading C:\Program Files\STNEXP\Queries\532185-1.str





chain nodes :

14 17

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 13

chain bonds : 1-14 5-17

ring bonds :

1-2 1-5 2-3 2-13 3-4 3-6 4-5 6-7 6-8 7-11 8-9 8-13 9-10 10-11

exact/norm bonds :

1-14 1-2 1-5 2-3 2-13 3-4 3-6 4-5 5-17 8-13

normalized bonds :

6-7 6-8 7-11 8-9 9-10 10-11

G1:Ir,Rh

G2:C1, Br, F, CH3, Et, n-Pr, i-Pr, Ph, MeO, EtO, n-PrO, i-PrO, CN, NO2

G3:O,S,N,Se

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 13:Atom 14:CLASS 17:CLASS

L1 STRUCTURE UPLOADED

=> d his

(FILE 'HOME' ENTERED AT 10:31:02 ON 22 JUN 2009)

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FILE 'REGISTRY' ENTERED AT 10:31:14 ON 22 JUN 2009
          STRUCTURE UPLOADED
```

=> d 11

L1

L1 HAS NO ANSWERS

L1STR

G1 Ir,Rh

G2 Cl,Br,F,Me,Et,n-Pr,i-Pr,Ph,MeO,EtO,n-PrO,i-PrO,CN,NO2

G3 O, S, N, Se

Structure attributes must be viewed using STN Express query preparation.

=> s 11

SAMPLE SEARCH INITIATED 10:31:44 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED -0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.10

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*

BATCH \*\*COMPLETE\*\* PROJECTED ITERATIONS: 0 TO

0 PROJECTED ANSWERS: 0 TO

L2 0 SEA SSS SAM L1

=> s 11 full

FULL SEARCH INITIATED 10:32:05 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED -0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

L3 0 SEA SSS FUL L1

=>

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Executing the logoff script...

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SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST 186.36 186.58

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